WEST Search History



DATE: Monday, May 14, 2007

Hide?	<u>Set</u> Name	Query	<u>Hit</u> Count
	DB=F	PGPB; PLUR=YES; OP=ADJ	
	L11	110 and (phosphonium salt or phosphonium compound or unsaturated phosphonium or olefinic phosphonium.CLM.)	1
	L10	19 and (acetal or aldehyde acetal or acetal aldehyde or acetal-containing aldehyde or dioxane or dioxalane.CLM.)	19
	L9	phytofluene or \$hexahydrolycopene.CLM.	124
	DB=F	PGPB, USPT, USOC, EPAB, JPAB, DWPI; PLUR=YES; OP=ADJ	
	L8	16 and (organic acid or acetic acid or formic acid or citric acid)	19
	L7	15 and (hydrolyz\$ or hydrolys\$ or cleav\$)	1
	L6	11 and 14	19
	L5	13 and 14	2
	L4	aldehyde acetal or acetal-containing aldehyde or dioxane or dioxalane	151816
	L3	11 and 12	5
	L2	phosphonium salt or phosphonium compound or unsaturated phosphonium compound or unsaturated phosphonium salt or olefinic phosphonium compound of olefinic phosphonium salt	18932
	L1	phytofluene or \$hexahydrolycopene	224

END OF SEARCH HISTORY

=> d his

(FILE 'HOME' ENTERED AT 15:24:33 ON 14 MAY 2007)

FILE 'REGISTRY' ENTERED AT 15:24:50 ON 14 MAY 2007

L1 STRUCTURE UPLOADED

L2 STRUCTURE UPLOADED

L3 0 S L2

L4 4 S L2 FULL

FILE 'HCAPLUS, CHEMCATS' ENTERED AT 15:26:50 ON 14 MAY 2007

L5 5 S L4

=> d his

L13

(FILE 'HOME' ENTERED AT 15:31:37 ON 14 MAY 2007)

	FILE	'REGISTRY'	ENTERED	AT	15:32:02	ON	14	MAY	2007
L1		STRU	CTURE UP	LOAI	DED				

2 S HENRICH, K/AU

L2 0 S L1

L3 0 S L1 FULL

FILE 'HCAPLUS, HCAOLD, USPATFULL, EPFULL' ENTERED AT 15:33:01 ON 14 MAY 2007

	2007		
L4	819	S	PHYTOFLUENE OR ?HEXAHYDROLYCOPENE
L5	72044	S	PHOSPHONIUM COMPOUND OR PHOSPHONIUM SALT OR UNSATURATED PHOSP
L6,	4	S	L4 AND L5
L7			L4 AND (ACETAL OR ALDEHYDE ACETAL OR ACETAL-CONTAINING ALDEHY
L8	20	S	L7 AND (CONDENS? OR COUPLE? OR COUPLING OR WITTIG?)
L9			L8 AND (HYDROLYS? OR HYDROLYZ? OR CLEAV?)
L10	19	S	L9 AND (ORGANIC ACID OR CARBOXYLIC ACID OR CITRIC ACID)
L11	187	S	ERNST, H/AU
L12	0	S	L11 AND PHYTOFLUENE

C:\Program Files\Stnexp\Queries\503-3.str

chain nodes:

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 18 19 20 21

ring nodes:

16 17 22 23 24 25

chain bonds:

1-2 2-3 2-18 3-4 4-5 5-6 6-7 6-19 7-8 8-9 9-10 10-11 10-20 11-12 12-13 13-14 14-15 14-21 15-16

ring bonds:

16-17 16-22 17-24 22-23 23-25 24-25

exact bonds:

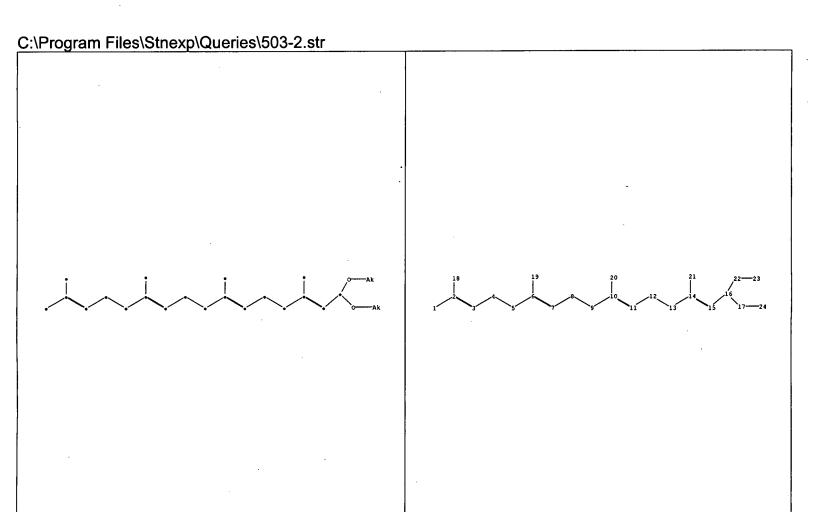
1-2 2-3 2-18 3-4 4-5 5-6 6-7 6-19 7-8 8-9 9-10 10-11 10-20 11-12 12-13 13-14 14-15 14-21 15-16 16-17 16-22 17-24 22-23 23-25 24-25

isolated ring systems:

containing 16:

Match level:

1:CLASS2:CLASS3:CLASS4:CLASS5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS 12:CLASS13:CLASS14:CLASS15:CLASS16:CLASS17:CLASS18:CLASS19:CLASS20:CLASS21:CLASS 22:CLASS23:Atom 24:Atom 25:Atom



chain nodes:

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 18 19 20 21 23 24

ring/chain nodes:

17 22

chain bonds:

1-2 2-3 2-18 3-4 4-5 5-6 6-7 6-19 7-8 8-9 9-10 10-11 10-20 11-12 12-13 13-14 14-15 14-21 15-16 16-17 16-22 17-24 22-23

exact/norm bonds:

16-17 16-22 17-24 22-23

exact bonds:

1-2 2-3 2-18 3-4 4-5 5-6 6-7 6-19 7-8 8-9 9-10 10-11 10-20 11-12 12-13 13-14 14-15 14-21 15-16

Match level:

1:CLASS2:CLASS3:CLASS4:CLASS5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS 12:CLASS13:CLASS14:CLASS15:CLASS16:CLASS17:CLASS18:CLASS19:CLASS20:CLASS21:CLASS 22:CLASS23:CLASS24:CLASS

Element Count:

Node 23: Limited

C.C1-9

Node 24: Limited

C,C1-9